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NEWS 6 JAN 28 USGENE now provides USPTO sequence data within 3 days of publication
NEWS 7 JAN 28 TOXCENTER enhanced with reloaded MEDLINE segment
NEWS 8 JAN 28 MEDLINE and LMEDLINE reloaded with enhancements
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NEWS 11 FEB 25 IFIREF reloaded with enhancements
NEWS 12 FEB 25 IMSPRODUCT reloaded with enhancements
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NEWS 16 MAR 31 CA/CAplus and CASREACT patent number format for U.S. applications updated
NEWS 17 MAR 31 LPCI now available as a replacement to LDPCI
NEWS 18 MAR 31 EMBASE, EMBAL, and LEMBASE reloaded with enhancements
NEWS 19 APR 04 STN AnaVist, Version 1, to be discontinued
NEWS 20 APR 15 WPIDS, WPINDEX, and WPIX enhanced with new predefined hit display formats
NEWS 21 APR 28 EMBASE Controlled Term thesaurus enhanced
NEWS 22 APR 28 IMSRESEARCH reloaded with enhancements
NEWS 23 MAY 30 INPAFAMDB now available on STN for patent family searching
NEWS 24 MAY 30 DGENE, PCTGEN, and USGENE enhanced with new homology sequence search option

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AND CURRENT DISCOVER FILE IS DATED 20 FEBRUARY 2008

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STRUCTURE FILE UPDATES: 3 JUN 2008 HIGHEST RN 1025148-33-7
DICTIONARY FILE UPDATES: 3 JUN 2008 HIGHEST RN 1025148-33-7

New CAS Information Use Policies, enter HELP USAGETERMS for details.

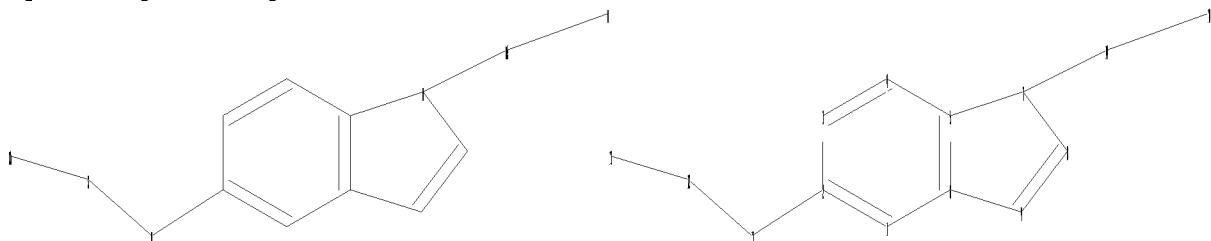
TSCA INFORMATION NOW CURRENT THROUGH January 9, 2008.

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REGISTRY includes numerically searchable data for experimental and predicted properties as well as tags indicating availability of experimental property data in the original document. For information on property searching in REGISTRY, refer to:

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=>
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10 11 12 13 14
ring nodes :
1 2 3 4 5 6 7 8 9
chain bonds :
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ring bonds :
1-2 1-6 2-3 3-4 4-5 5-6 5-7 6-9 7-8 8-9
exact/norm bonds :
2-11 5-7 6-9 7-8 7-14 8-9 10-14 11-12

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exact bonds :
12-13
normalized bonds :
1-2 1-6 2-3 3-4 4-5 5-6

Match level :
1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:Atom 9:Atom 10:CLASS
11:CLASS 12:CLASS 13:Atom 14:CLASS

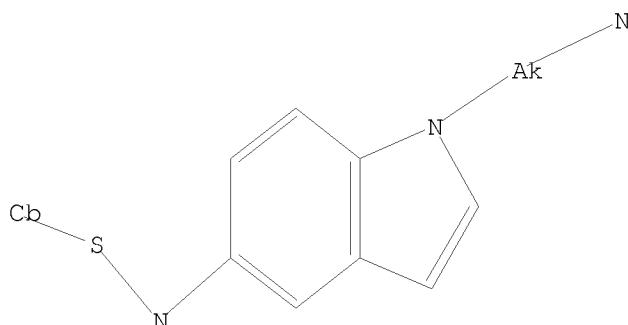
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L1 STRUCTURE UPLOADED

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L1 HAS NO ANSWERS
L1           STR

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Structure attributes must be viewed using STN Express query preparation.

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SAMPLE SEARCH INITIATED 14:55:18 FILE 'REGISTRY'
SAMPLE SCREEN SEARCH COMPLETED -       1091 TO ITERATE

100.0% PROCESSED       1091 ITERATIONS                           2 ANSWERS
SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE    **COMPLETE**
                          BATCH    **COMPLETE**
PROJECTED ITERATIONS:       19839 TO    23801
PROJECTED ANSWERS:          2 TO      124

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L2 2 SEA SSS SAM L1

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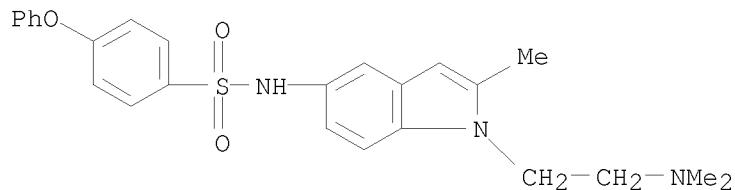
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L2    2 ANSWERS    REGISTRY   COPYRIGHT 2008 ACS on STN
IN    Benzenesulfonamide, N-[1-[2-(dimethylamino)ethyl]-2-methyl-1H-indol-5-yl]-
      4-phenoxy-
MF    C25 H27 N3 O3 S

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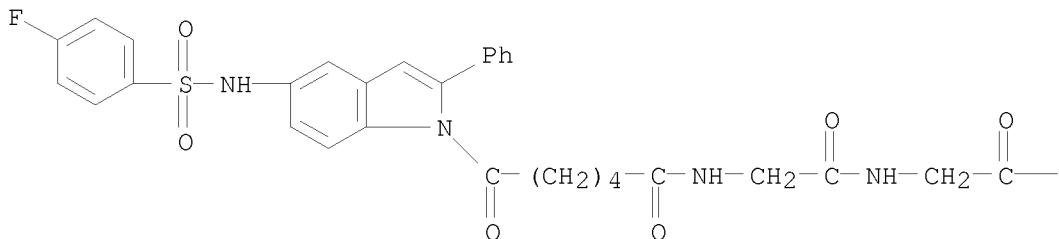


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):1

L2 2 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN
 IN Glycinamide, N-[6-[5-[(4-fluorophenyl)sulfonyl]amino]-2-phenyl-1H-indol-1-yl]-1,6-dioxohexylglycyl-
 MF C30 H30 F N5 O6 S

PAGE 1-A



PAGE 1-B

— NH₂

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

ALL ANSWERS HAVE BEEN SCANNED

=> s l1 sss full
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 FULL SCREEN SEARCH COMPLETED - 21961 TO ITERATE

100.0% PROCESSED 21961 ITERATIONS
 SEARCH TIME: 00.00.01

36 ANSWERS

L3 36 SEA SSS FUL L1

=> file caplus

COST IN U.S. DOLLARS	SINCE FILE ENTRY	TOTAL SESSION
FULL ESTIMATED COST	179.28	179.49

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 FILE LAST UPDATED: 3 Jun 2008 (20080603/ED)

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=> s 13
 L4 6 L3

=> d ibib abs hitstr 6

L4 ANSWER 6 OF 6 CAPLUS COPYRIGHT 2008 ACS on STN
 ACCESSION NUMBER: 2004:725572 CAPLUS
 DOCUMENT NUMBER: 142:211383
 TITLE: Medicinal Chemistry Driven Approaches Toward Novel and Selective Serotonin 5-HT₆ Receptor Ligands
 Holenz, Joerg; Merce, Ramon; Diaz, Jose Luis; Guitart, Xavier; Codony, Xavier; Dordal, Alberto; Romero, Gonzalo; Torrens, Antoni; Mas, Josep; Andaluz, Blas; Hernandez, Susana; Monroy, Xavier; Sanchez, Elisabeth; Hernandez, Enrique; Perez, Raquel; Cubi, Roger; Sanfeliu, Olga; Buschmann, Helmut
 AUTHOR(S):
 CORPORATE SOURCE: Departments of Medicinal Chemistry, Discovery Biology and Discovery Chemistry, Laboratorios Dr. Esteve S.A., Barcelona, 08041, Spain
 SOURCE: Journal of Medicinal Chemistry (2005), 48(6), 1781-1795
 CODEN: JMCMAR; ISSN: 0022-2623
 PUBLISHER: American Chemical Society
 DOCUMENT TYPE: Journal
 LANGUAGE: English
 OTHER SOURCE(S): CASREACT 142:211383
 AB Based on a medicinal chemical guided hypothetical pharmacophore model, novel series of indolyl sulfonamides have been designed and prepared as selective and high-affinity serotonin 5-HT₆ receptor ligands. Furthermore, based on a screening approach of a discovery library, a series of benzoxazineperidinyl sulfonamides were identified as selective 5-HT₆ ligands. Many of the compds. described in this paper possess excellent affinities, displaying pKi values greater than 8 (some even >9) and high selectivities against a wide range (>50) of other CNS relevant receptors.

First, structure-affinity relationships of these ligands are discussed. In terms of functionality, high-affinity antagonists, as well as agonists and even partial agonists, were prepared. Compds. 19c and 19g represent the highest-affinity 5-HT₆ agonists ever reported in the literature. These valuable tool compds. should allow for the detailed study of the role of the 5-HT₆ receptor in relevant animal models of disorders such as cognition deficits, depression, anxiety, or obesity.

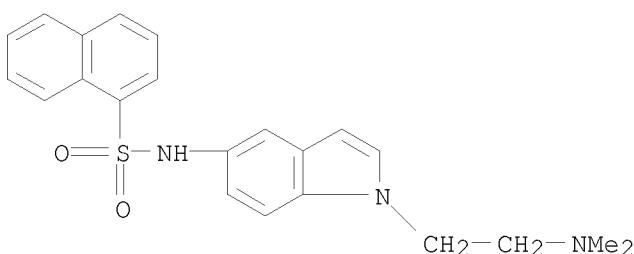
IT 753020-75-6P 753020-76-7P 753020-79-0P
 753020-80-3P 753020-82-5P 753020-83-6P
 753020-84-7P

RL: DMA (Drug mechanism of action); PAC (Pharmacological activity); PRP (Properties); SPN (Synthetic preparation); BIOL (Biological study); PREP (Preparation)

(medicinal chemical driven approaches toward novel and selective serotonin 5-HT₆ receptor ligands)

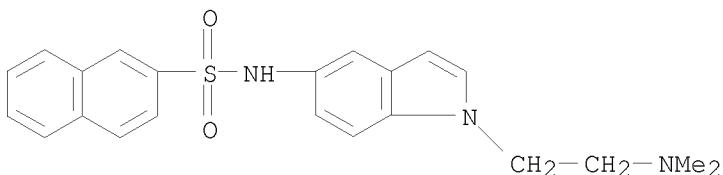
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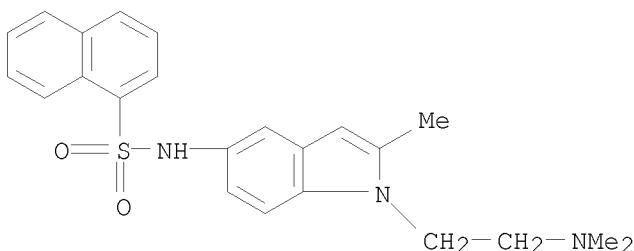
RN 753020-76-7 CAPLUS

CN 2-Naphthalenesulfonamide, N-[1-[2-(dimethylamino)ethyl]-1H-indol-5-yl]-(CA INDEX NAME)



RN 753020-79-0 CAPLUS

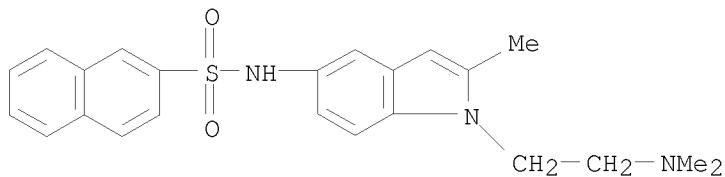
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RN 753020-80-3 CAPLUS

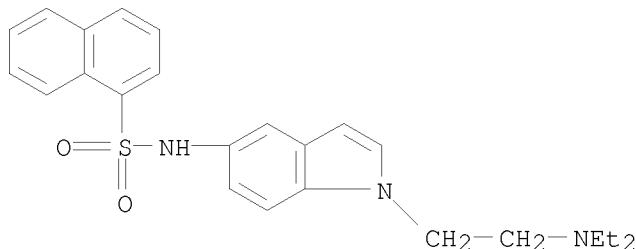
CN 2-Naphthalenesulfonamide, N-[1-[2-(dimethylamino)ethyl]-2-methyl-1H-indol-

5-y1]- (CA INDEX NAME)



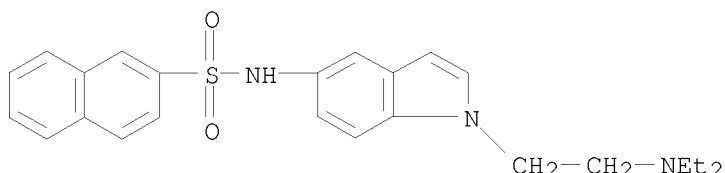
RN 753020-82-5 CAPLUS

CN 1-Naphthalenesulfonamide, N-[1-[2-(diethylamino)ethyl]-1H-indol-5-yl]-
(CA INDEX NAME)



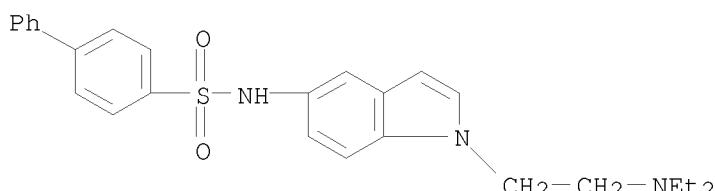
RN 753020-83-6 CAPLUS

CN 2-Naphthalenesulfonamide, N-[1-[2-(diethylamino)ethyl]-1H-indol-5-yl]-
(CA INDEX NAME)



RN 753020-84-7 CAPLUS

CN [1,1'-Biphenyl]-4-sulfonamide, N-[1-[2-(diethylamino)ethyl]-1H-indol-5-yl]-
(CA INDEX NAME)



REFERENCE COUNT:

68

THERE ARE 68 CITED REFERENCES AVAILABLE FOR THIS
RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

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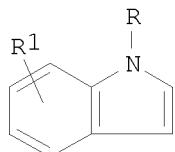
L4 ANSWER 1 OF 6 CAPLUS COPYRIGHT 2008 ACS on STN
ACCESSION NUMBER: 2007:197836 CAPLUS

DOCUMENT NUMBER: 146:252104
 TITLE: Preparation of substituted indoles and their use as PAI-1 inhibitors
 INVENTOR(S): Hu, Baihua; Jetter, James W.
 PATENT ASSIGNEE(S): Wyeth, John, and Brother Ltd., USA
 SOURCE: PCT Int. Appl., 54pp.
 CODEN: PIXXD2
 DOCUMENT TYPE: Patent
 LANGUAGE: English
 FAMILY ACC. NUM. COUNT: 1
 PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
WO 2007022321	A2	20070222	WO 2006-US32066	20060816
WO 2007022321	A3	20070510		
W: AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BW, BY, BZ, CA, CH, CN, CO, CR, CU, CZ, DE, DK, DM, DZ, EC, EE, EG, ES, FI, GB, GD, GE, GH, GM, HN, HR, ID, IL, IN, IS, JP, KE, KG, KM, KN, KP, KR, KZ, LA, LC, LK, LR, LS, LT, LU, LV, LY, MA, MD, MG, MK, MN, MW, MX, MY, MZ, NA, NG, NI, NO, NZ, OM, PG, PH, PL, PT, RO, RS, RU, SC, SD, SE, SG, SK, SL, SM, SY, TJ, TM, TN, TR, TT, TZ, UA, UG, US, UZ, VC, VN, ZA, ZM, ZW				
RW: AT, BE, BG, CH, CY, CZ, DE, DK, EE, ES, FI, FR, GB, GR, HU, IE, IS, IT, LT, LU, LV, MC, NL, PL, PT, RO, SE, SI, SK, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE, SN, TD, TG, BW, GH, GM, KE, LS, MW, MZ, NA, SD, SL, SZ, TZ, UG, ZM, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM, AP, EA, EP, OA				
AU 2006279496	A1	20070222	AU 2006-279496	20060816
CA 2617372	A1	20070222	CA 2006-2617372	20060816
US 20070043101	A1	20070222	US 2006-505527	20060816
EP 1919866	A2	20080514	EP 2006-801683	20060816
R: AT, BE, BG, CH, CY, CZ, DE, DK, EE, ES, FI, FR, GB, GR, HU, IE, IS, IT, LI, LT, LU, LV, MC, NL, PL, PT, RO, SE, SI, SK, TR				
PRIORITY APPLN. INFO.:			US 2005-708834P	P 20050817
			WO 2006-US32066	W 20060816

OTHER SOURCE(S): MARPAT 146:252104

GI



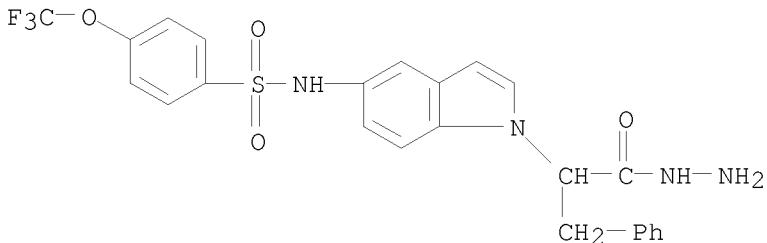
AB The invention relates to indole derivs. I [R is p-R₂C₆H₄(CH₂)₁₋₄, where R₂ is alkyl, and R₁ is a sulfonylamino or ureido group; or R is R₃C₆H₄(CH₂)₀₋₄CHR₄, where R₃ is H, a carboxyalkoxy, carbamoyl, or carbonyl-amino acid group and R₄ is H, CO₂H, or CONHNH₂ and R₁ is a sulfonylamino group; or R is R₅CO(CH₂)₁₋₄, where R₅ is OH, alkoxy, or an amino acid residue and R₁ is a sulfonylamino group] for use as PAI-1 inhibitors. Thus, N-[[[1-(4-tert-butylbenzyl)-1H-indol-5-yl]amino]carbonyl]-L-phenylalanine was prepared by treating 1-(4-tert-butylbenzyl)-1H-indol-5-amine (preparation given) with 2-isocyanato-3-phenylpropionic acid Et ester.

IT 926024-84-2P

RL: PAC (Pharmacological activity); RCT (Reactant); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); RACT (Reactant or reagent); USES (Uses)
(preparation of substituted indoles and their use as PAI-1 inhibitors)

RN 926024-84-2 CAPLUS

CN 1H-Indole-1-acetic acid, α -(phenylmethyl)-5-[[(4-(trifluoromethoxy)phenyl)sulfonyl]amino]-, hydrazide (CA INDEX NAME)



IT 926024-62-6P 926024-64-8P 926024-66-0P

926024-68-2P 926024-70-6P 926024-74-0P

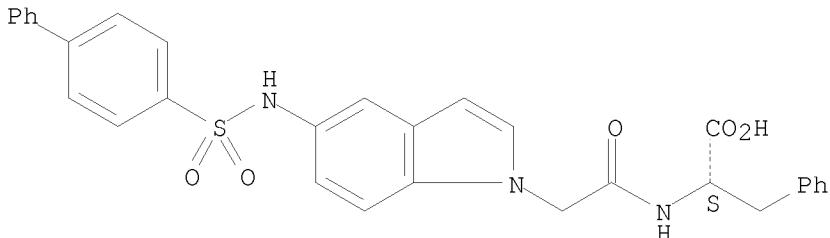
RL: PAC (Pharmacological activity); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)

(preparation of substituted indoles and their use as PAI-1 inhibitors)

RN 926024-62-6 CAPLUS

CN L-Phenylalanine, N-[2-[5-[[(1,1'-biphenyl)-4-ylsulfonyl]amino]-1H-indol-1-yl]acetyl]- (CA INDEX NAME)

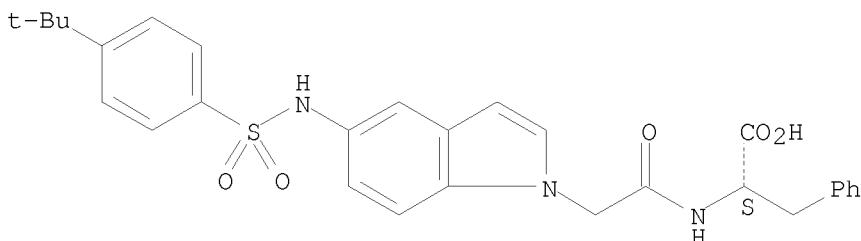
Absolute stereochemistry.



RN 926024-64-8 CAPLUS

CN L-Phenylalanine, N-[2-[5-[[(4-(1,1-dimethylethyl)phenyl)sulfonyl]amino]-1H-indol-1-yl]acetyl]- (CA INDEX NAME)

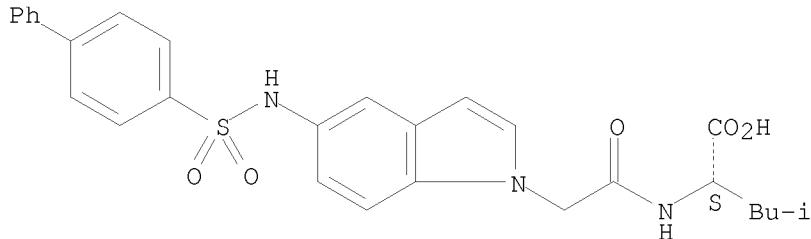
Absolute stereochemistry.



RN 926024-66-0 CAPLUS

CN L-Leucine, N-[2-[5-((1,1'-biphenyl)-4-ylsulfonyl)amino]-1H-indol-1-yl]acetyl]- (CA INDEX NAME)

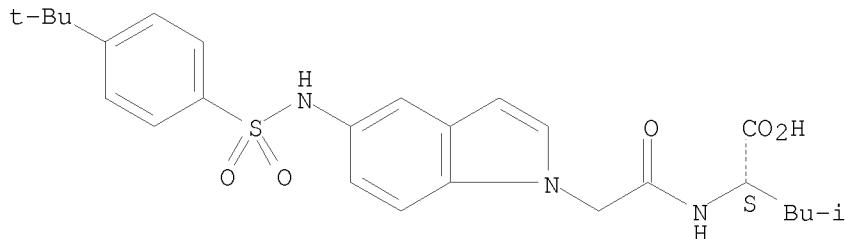
Absolute stereochemistry.



RN 926024-68-2 CAPLUS

CN L-Leucine, N-[2-[5-[[4-(1,1-dimethylethyl)phenyl]sulfonyl]amino]-1H-indol-1-yl]acetyl]- (CA INDEX NAME)

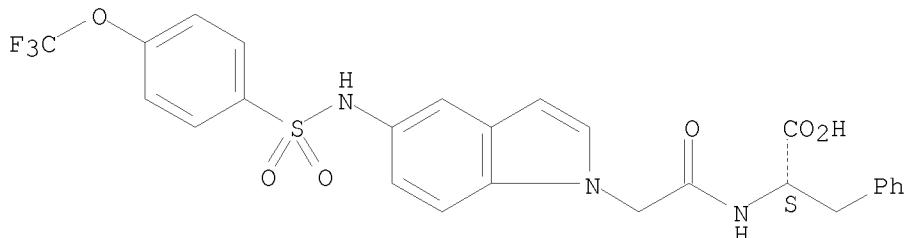
Absolute stereochemistry.



RN 926024-70-6 CAPLUS

CN L-Phenylalanine, N-[2-[5-[[4-(trifluoromethoxy)phenyl]sulfonyl]amino]-1H-indol-1-yl]acetyl]- (CA INDEX NAME)

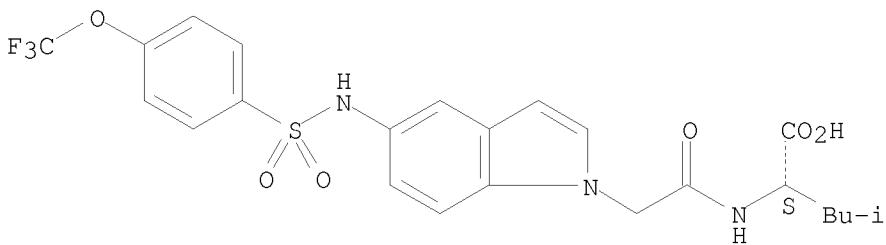
Absolute stereochemistry.



RN 926024-74-0 CAPLUS

CN L-Leucine, N-[2-[5-[[4-(trifluoromethoxy)phenyl]sulfonyl]amino]-1H-indol-1-yl]acetyl]- (CA INDEX NAME)

Absolute stereochemistry.



L4 ANSWER 2 OF 6 CAPLUS COPYRIGHT 2008 ACS on STN

ACCESSION NUMBER: 2006:1244611 CAPLUS

DOCUMENT NUMBER: 146:142447

TITLE: An engineered linker capable of promoting on-resin reactions for microwave-assisted solid-phase organic synthesis

AUTHOR(S): Sun, Li-Ping; Dai, Wei-Min

CORPORATE SOURCE: Department of Chemistry, The Hong Kong University of Science and Technology, Clear Water Bay, Kowloon, Hong Kong

SOURCE: Angewandte Chemie, International Edition (2006), 45(43), 7255-7258

PUBLISHER: Wiley-VCH Verlag GmbH & Co. KGaA

DOCUMENT TYPE: Journal

LANGUAGE: English

OTHER SOURCE(S): CASREACT 146:142447

AB A diglycine-containing linker was fabricated on Rink amide resin for dual functions: a) attachment of a scaffold and b) capture of metal ions for promoting on-resin reactions. The metal-catching feature of the linker proves essential for the solid-phase synthesis of indoles through microwave-assisted CuII-mediated heteroannulation.

IT 919490-32-7P 919490-34-9P 919490-36-1P

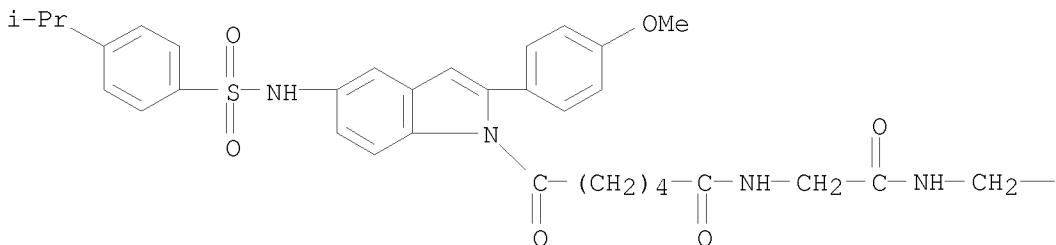
RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)

(microwave-assisted copper-mediated solid-phase synthesis of (arylsulfonylamino)indoles using diglycine-containing linker capable of catching metal ions and promoting on-resin reactions)

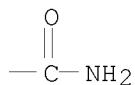
RN 919490-32-7 CAPLUS

CN Glycinamide, N-[6-[2-(4-methoxyphenyl)-5-[[[4-(1-methylethyl)phenyl]sulfonyl]amino]-1H-indol-1-yl]-1,6-dioxohexyl]glycyl- (CA INDEX NAME)

PAGE 1-A

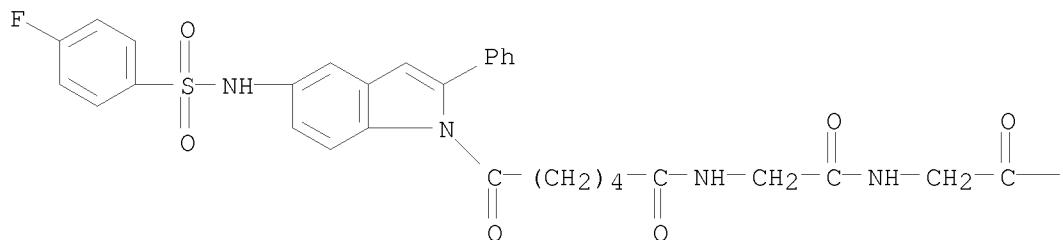


PAGE 1-B



RN 919490-34-9 CAPLUS
CN Glycinamide, N-[6-[5-[(4-fluorophenyl)sulfonyl]amino]-2-phenyl-1H-indol-1-yl]-1,6-dioxohexylglycyl- (CA INDEX NAME)

PAGE 1-A

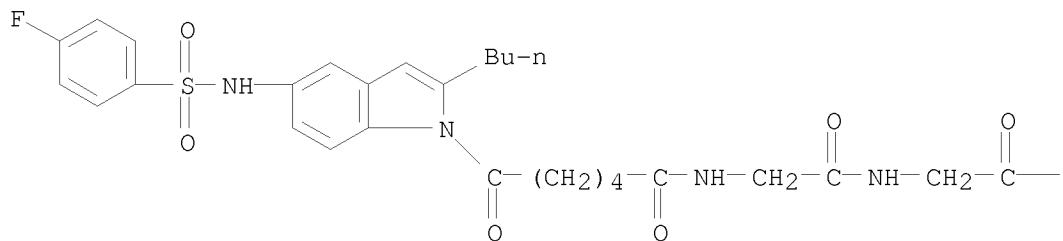


PAGE 1-B



RN 919490-36-1 CAPLUS
CN Glycinamide, N-[6-[2-butyl-5-[(4-fluorophenyl)sulfonyl]amino]-1H-indol-1-yl]-1,6-dioxohexylglycyl- (CA INDEX NAME)

PAGE 1-A



PAGE 1-B

— NH₂

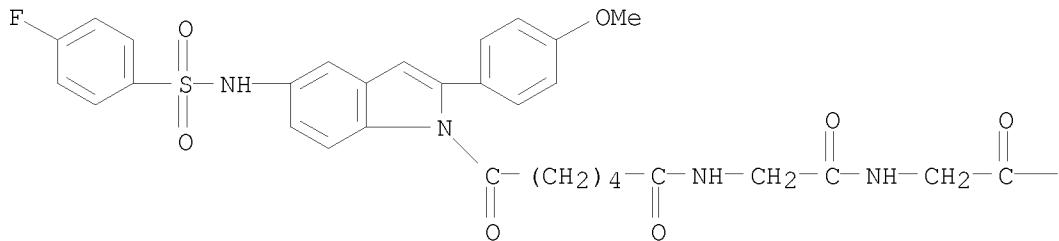
IT 919490-25-8P 919490-27-0P 919490-28-1P
919490-30-5P

RL: SPN (Synthetic preparation); PREP (Preparation)
(microwave-assisted copper-mediated solid-phase synthesis of
(arylsulfonylamino)indoles using diglycine-containing linker capable of
catching metal ions and promoting on-resin reactions)

RN 919490-25-8 CAPLUS

CN Glycinamide, N-[6-[5-[(4-fluorophenyl)sulfonyl]amino]-2-(4-methoxyphenyl)-
1H-indol-1-yl]-1,6-dioxohexyl]glycyl- (CA INDEX NAME)

PAGE 1-A

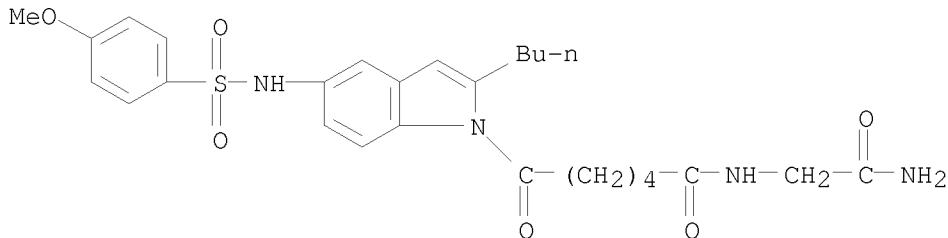


PAGE 1-B

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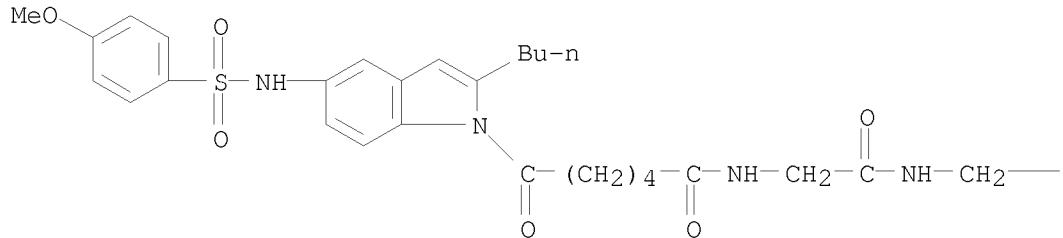
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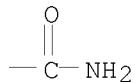
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CN Glycinamide, N-[6-[2-butyl-5-[(4-methoxyphenyl)sulfonyl]amino]-1H-indol-1-yl]-1,6-dioxohexylglycyl- (CA INDEX NAME)

PAGE 1-A



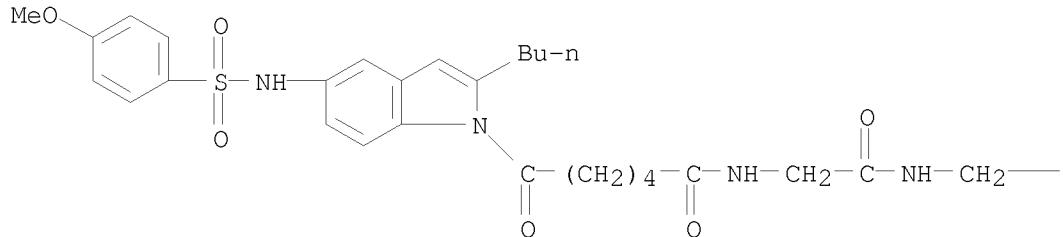
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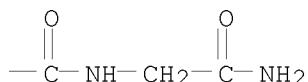
RN 919490-30-5 CAPLUS

CN Glycinamide, N-[6-[2-butyl-5-[(4-methoxyphenyl)sulfonyl]amino]-1H-indol-1-yl]-1,6-dioxohexylglycylglycyl- (CA INDEX NAME)

PAGE 1-A



PAGE 1-B



REFERENCE COUNT:

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THERE ARE 5 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L4 ANSWER 3 OF 6 CAPLUS COPYRIGHT 2008 ACS on STN

ACCESSION NUMBER: 2005:136598 CAPLUS

DOCUMENT NUMBER: 142:240323

TITLE: Active substance combination comprising a compound

with NPY receptor affinity and a compound with 5-HT6
 receptor affinity
 INVENTOR(S):
 Torrens Jover, Antoni; Mas Prio, Josep; Dordal Zueras,
 Alberto; Codony Soler, Xavier; Merce Vidal, Ramon;
 Aurelio Castrillo Perez, Jose; Frigola Constansa,
 Jordi; Buschmann, Helmut-Heinrich
 PATENT ASSIGNEE(S):
 Laboratorios del Esteve S. A., Spain
 SOURCE:
 PCT Int. Appl., 427 pp.
 CODEN: PIXXD2
 DOCUMENT TYPE:
 Patent
 LANGUAGE:
 English
 FAMILY ACC. NUM. COUNT: 1
 PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
WO 2005014045	A1	20050217	WO 2004-EP8514	20040729
W: AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BW, BY, BZ, CA, CH, CN, CO, CR, CU, CZ, DE, DK, DM, DZ, EC, EE, EG, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NA, NI, NO, NZ, OM, PG, PH, PL, PT, RO, RU, SC, SD, SE, SG, SK, SL, SY, TJ, TM, TN, TR, TT, TZ, UA, UG, US, UZ, VC, VN, YU, ZA, ZM, ZW				
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PRIORITY APPLN. INFO.:			ES 2003-1815	A 20030730
			WO 2004-EP8514	W 20040729

OTHER SOURCE(S): CASREACT 142:240323; MARPAT 142:240323
 GI

* STRUCTURE DIAGRAM TOO LARGE FOR DISPLAY - AVAILABLE VIA OFFLINE PRINT *

AB The present invention relates to an active substance combination comprising at least one compound I [R1-R4 = H, halo, alkyl, etc.; R5 = H, alkyl, (un)saturated cycloalkyl; R6-R9 = H, alkyl, (un)saturated cycloalkyl, etc.]; A = CHR18, CHR18CH2; B = alkyl, (un)saturated cycloalkyl, etc.; R10 = H, alkyl, (un)saturated cycloalkyl, etc.; R11 = alkyl, (un)saturated cycloalkyl, etc.; NR10R11 = (un)saturated heterocyclyl; R18 = H, alkyl, (un)saturated cycloalkyl, etc.] with neuropeptide Y-receptor affinity, preferably neuropeptide Y5-receptor affinity, and at least one compound with 5-HT6 receptor affinity (such as II [R1 = H, alkyl, Ph, CH2PH; R2 = NR4R5, (un)saturated (hetero)cycloalkyl, etc.; R3 = H, alkyl; R4, R5 = H, alkyl; or NR4R5 = (un)saturated heterocyclyl; A = (un)substituted (hetero)aryl; n = 0-4]], a medicament comprising said active substance combination, and the use of said active substance combination for the manufacture of a medicament. Synthesis of amides I and sulfonamides such as II is described in

examples. E.g., a multi-step synthesis of III.HCl, starting from 1-(tert-butoxycarbonyl)-4-piperidinone and Me anthranilate, was given. The amides I and sulfonamides such as II were tested against neuropeptide Y5 and 5-HT6 binding (data given for representative compds.).

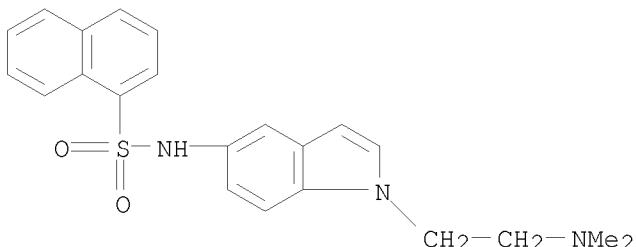
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RL: PAC (Pharmacological activity); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)

(preparation of amides and sulfonamides as components of active combination with NPY receptor affinity and 5-HT6 receptor affinity)

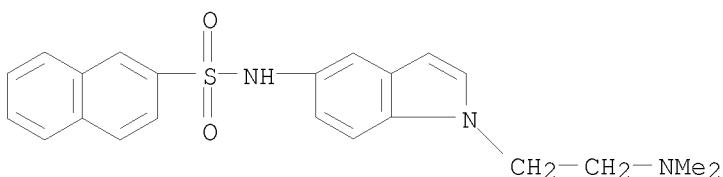
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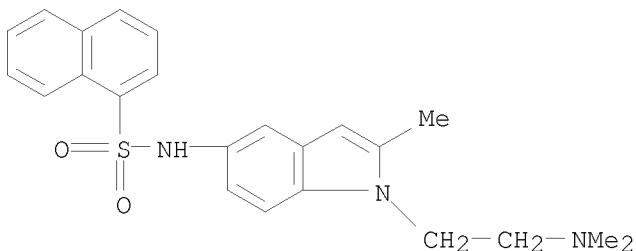
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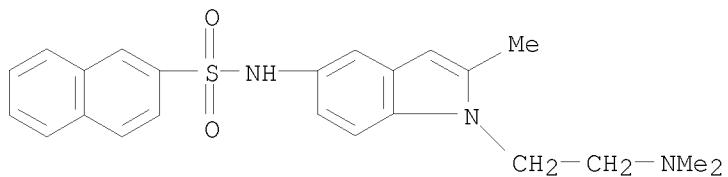


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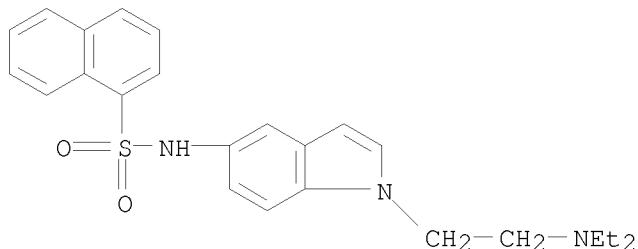
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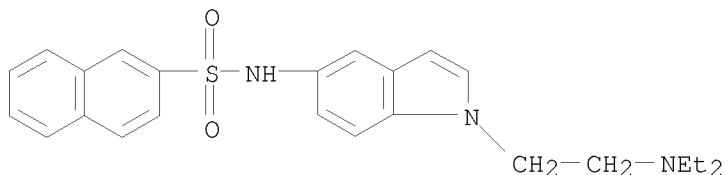
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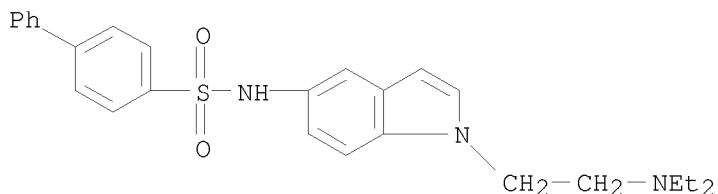
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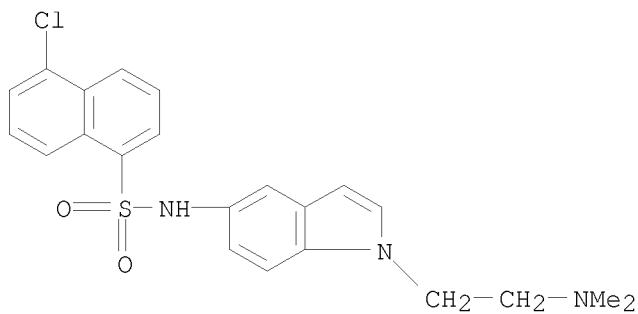
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CN 2-Naphthalenesulfonamide, N-[1-[2-(diethylamino)ethyl]-1H-indol-5-yl]- (CA INDEX NAME)



RN 753020-84-7 CAPLUS
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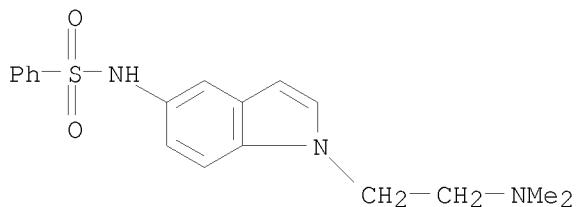


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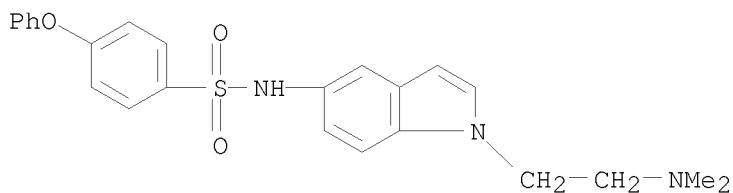
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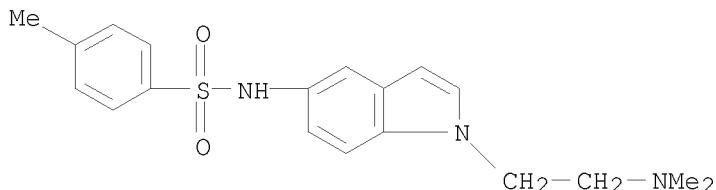
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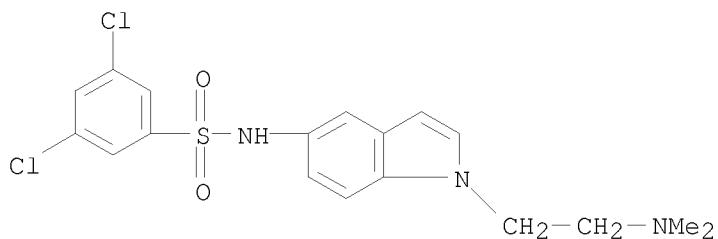
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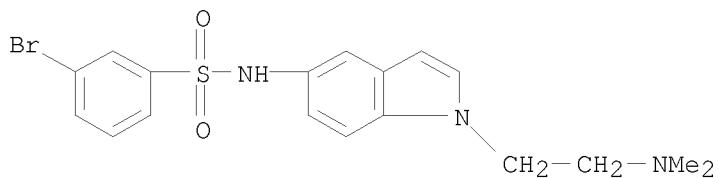
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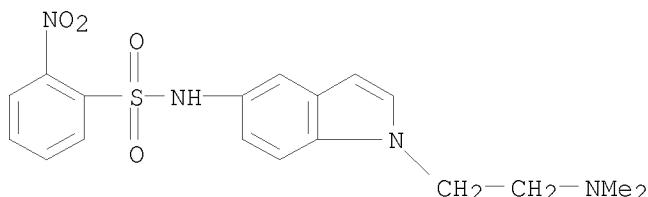
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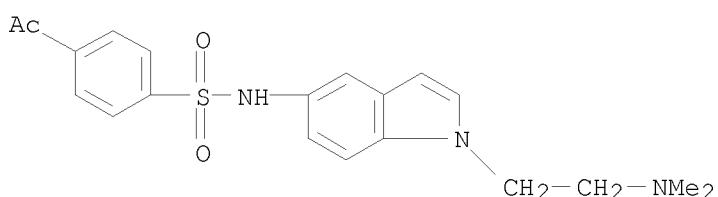
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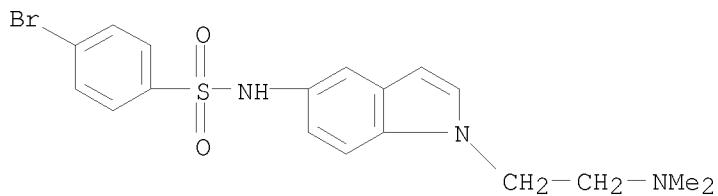
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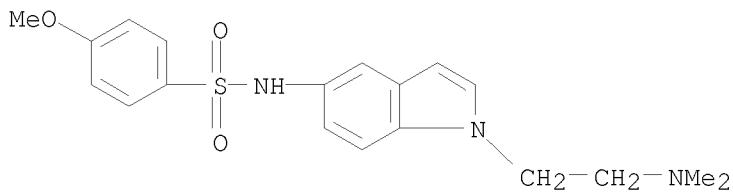
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(CA INDEX NAME)



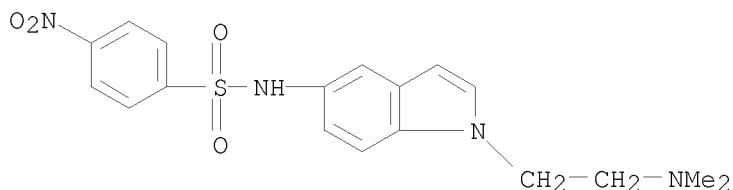
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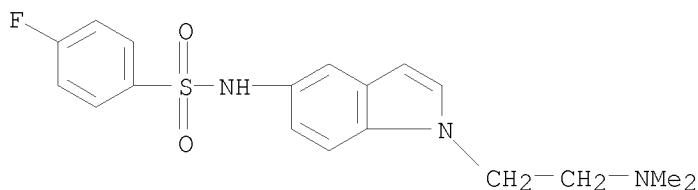
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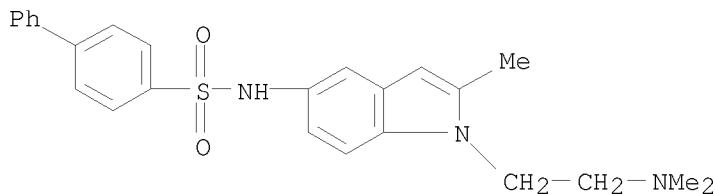
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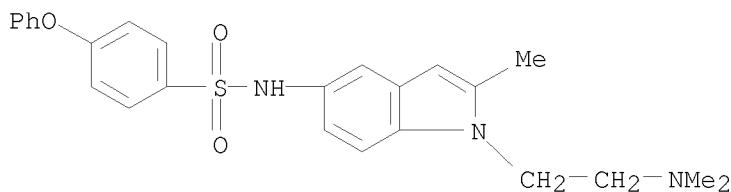
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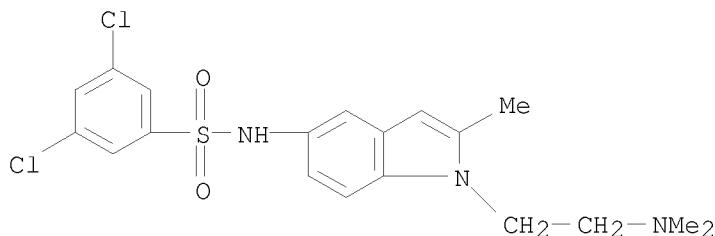
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CN Benzenesulfonamide, N-[1-[2-(dimethylamino)ethyl]-2-methyl-1H-indol-5-yl]-4-phenoxy- (CA INDEX NAME)



RN 844832-10-6 CAPLUS

CN Benzenesulfonamide, 3,5-dichloro-N-[1-[2-(dimethylamino)ethyl]-2-methyl-1H-indol-5-yl]- (CA INDEX NAME)



REFERENCE COUNT: 4 THERE ARE 4 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L4 ANSWER 4 OF 6 CAPLUS COPYRIGHT 2008 ACS on STN

ACCESSION NUMBER: 2005:136568 CAPLUS

DOCUMENT NUMBER: 142:240322

TITLE: Active substance combination comprising a compound with NPY receptor affinity and a compound with 5-HT6 receptor affinity

INVENTOR(S): Torrens Jover, Antoni; Mas Prio, Josep; Dordal Zueras, Alberto; Codony Soler, Xavier; Merce Vidal, Ramon; Aurelio Castrillo Perez, Jose; Frigola Constansa, Jordi; Buschmann, Helmut-Heinrich

PATENT ASSIGNEE(S): Laboratorios del Esteve S. A., Spain

SOURCE: PCT Int. Appl., 451 pp.

CODEN: PIXXD2

DOCUMENT TYPE: Patent

LANGUAGE: English

FAMILY ACC. NUM. COUNT: 1

PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
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WO 2005014000	A1	20050217	WO 2004-EP8515	20040729
W: AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BW, BY, BZ, CA, CH, CN, CO, CR, CU, CZ, DE, DK, DM, DZ, EC, EE, EG, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NA, NI, NO, NZ, OM, PG, PH, PL, PT, RO, RU, SC, SD, SE, SG, SK, SL, SY, TJ, TM, TN, TR, TT, TZ, UA, UG, US, UZ, VC, VN, YU, ZA, ZM, ZW				
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PRIORITY APPLN. INFO.:			ES 2003-1814	A 20030730
			WO 2004-EP8515	W 20040729

OTHER SOURCE(S): MARPAT 142:240322
GI

* STRUCTURE DIAGRAM TOO LARGE FOR DISPLAY - AVAILABLE VIA OFFLINE PRINT *

AB The present invention relates to an active substance combination comprising at least one compound I [R1-R4 = H, halo, alkyl, etc.; R5 = H, alkyl, (un)saturated (hetero)cycloalkyl; R6-R9 = H, alkyl, (un)saturated (hetero)cycloalkyl, etc.; A = CHR18, CHR18CH2; R10 = H, alkyl, (un)saturated cycloalkyl, etc.; R11 = alkyl, (un)saturated cycloalkyl, etc.; NR10R11 = (un)saturated heterocyclyl; R18 = H, alkyl, (un)saturated cycloalkyl, etc.]

with

neuropeptide Y-receptor affinity, preferably neuropeptide Y5-receptor affinity, and at least one compound with 5-HT6 receptor affinity (such as II [R1 = H, alkyl, Ph, CH2PH; R2 = NR4R5, (un)saturated (hetero)cycloalkyl, etc.; R3 = H, alkyl; R4, R5 = H, alkyl; or NR4R5 = (un)saturated heterocyclyl; A = (un)substituted (hetero)aryl; n = 0-4]), a medicament comprising said active substance combination, and the use of said active substance combination for the manufacture of a medicament. Synthesis of amides I and sulfonamides such as II is described in examples. Thus, reacting 6-chloro-1-(4-piperidinyl)-1,4-dihydro-2H-3,1-benzoxazinone hydrochloride with 2-(2-chloroacetamide)-2',5-dichlorobenzophenone in the presence of K2CO3 in DMF followed by treating of the free base with HCl/EtOH afforded 61% III.HCl. The amides I and sulfonamides such as II were tested against neuropeptide Y5 and 5-HT6 binding (data given for representative compds.).

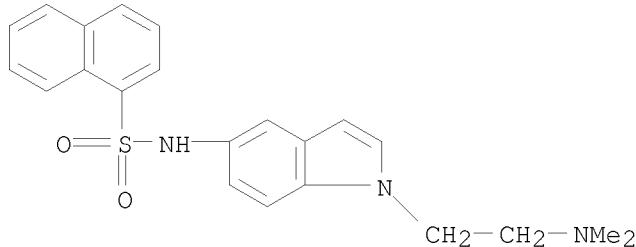
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844832-10-6P

RL: PAC (Pharmacological activity); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)

(preparation of amides and sulfonamides as components of active combination with NPY receptor affinity and 5-HT₆ receptor affinity)

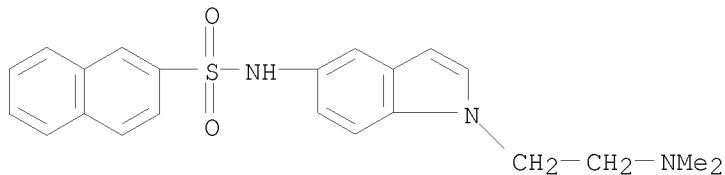
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(CA INDEX NAME)



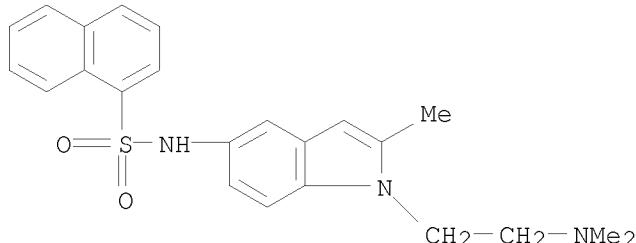
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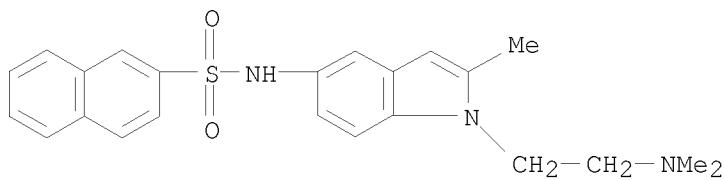
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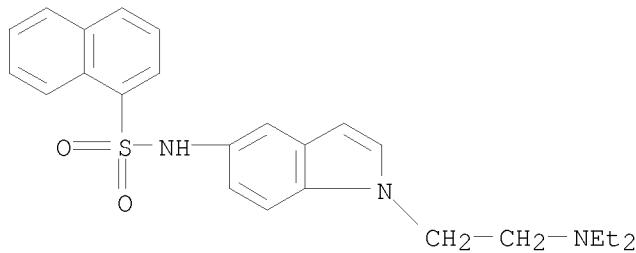
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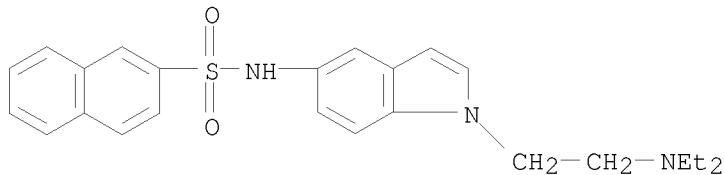
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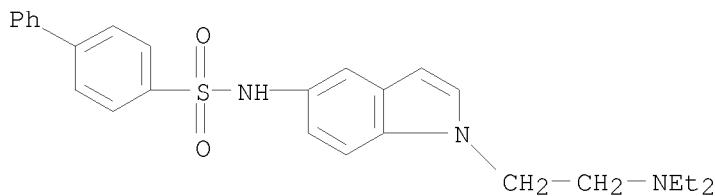
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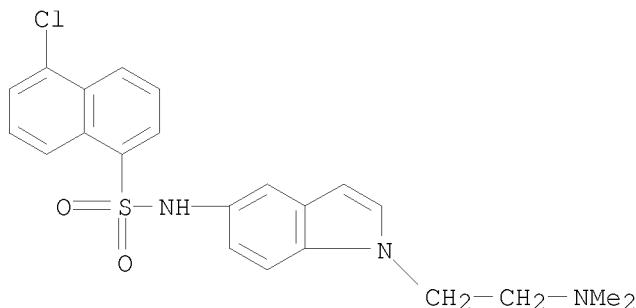
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(CA INDEX NAME)



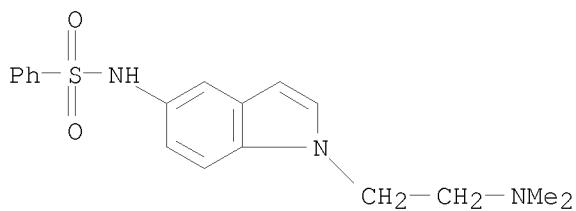
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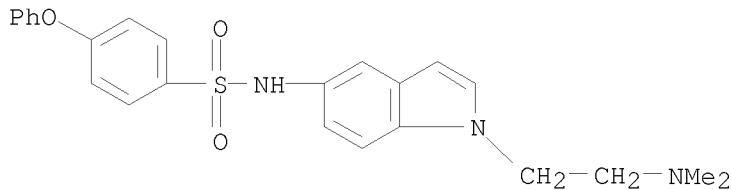


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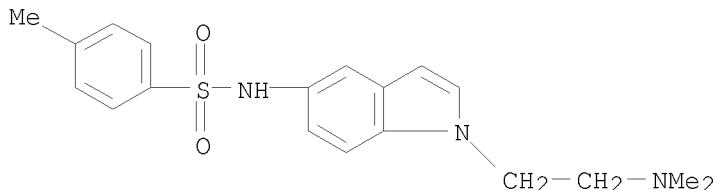
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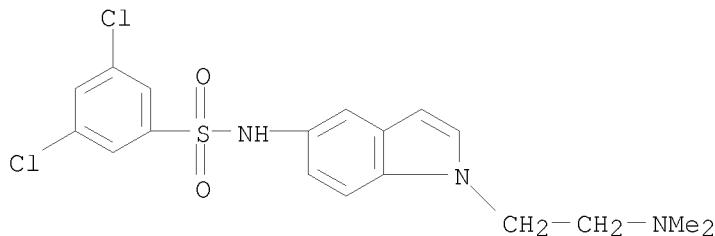
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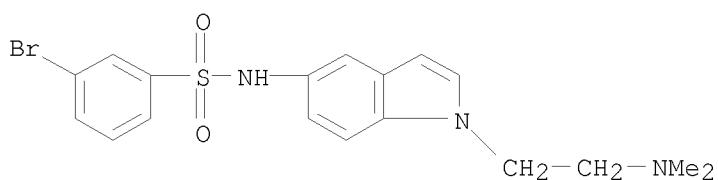
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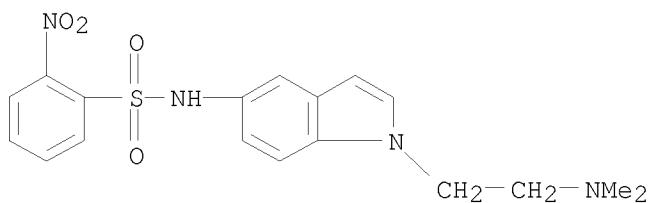
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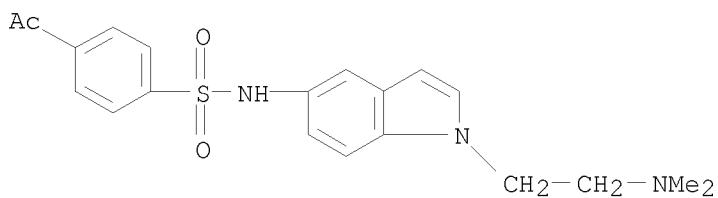
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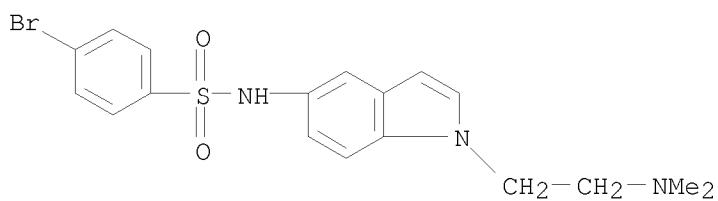
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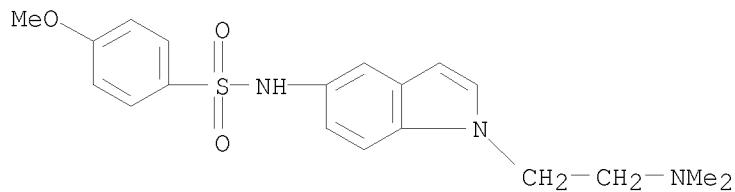
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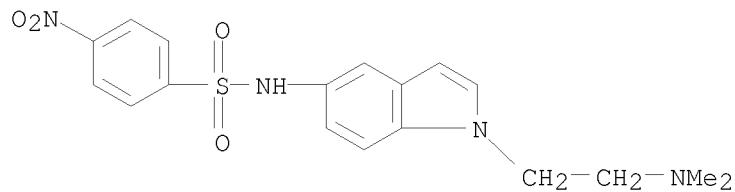


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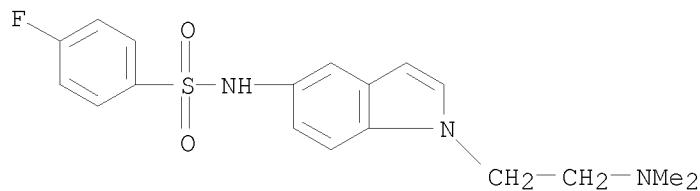
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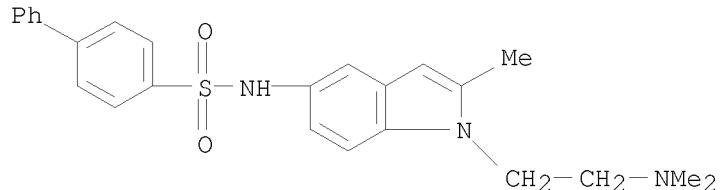
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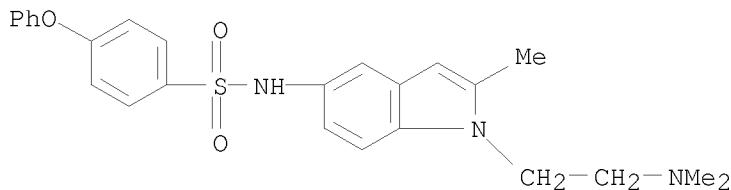
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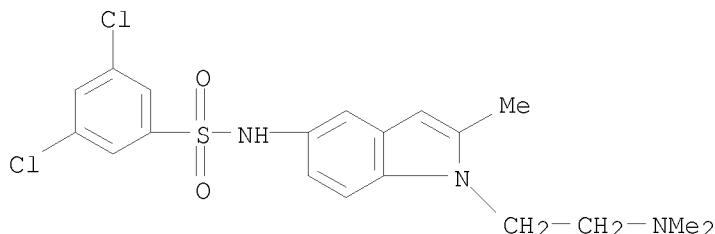
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RN 844832-10-6 CAPLUS

CN Benzenesulfonamide, 3,5-dichloro-N-[1-[2-(dimethylamino)ethyl]-2-methyl-1H-indol-5-yl]- (CA INDEX NAME)



REFERENCE COUNT: 5 THERE ARE 5 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L4 ANSWER 5 OF 6 CAPLUS COPYRIGHT 2008 ACS on STN

ACCESSION NUMBER: 2005:136549 CAPLUS

DOCUMENT NUMBER: 142:240310

TITLE: Preparation of indol-5-yl sulfonamide derivatives and their use as 5-HT6 modulators

INVENTOR(S): Merce Vidal, Ramon; Codony Soler, Xavier; Dordal Zueras, Alberto

PATENT ASSIGNEE(S): Laboratorios del Esteve S. A., Spain

SOURCE: PCT Int. Appl., 123 pp.

CODEN: PIXXD2

DOCUMENT TYPE: Patent

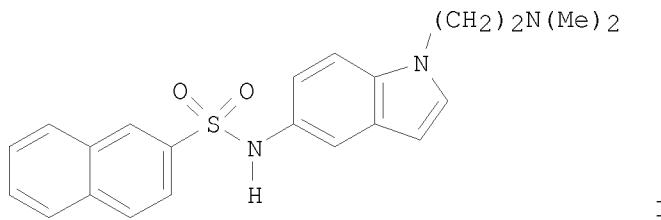
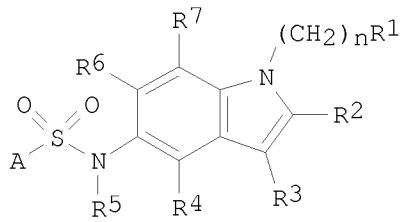
LANGUAGE: English

FAMILY ACC. NUM. COUNT: 1

PATENT INFORMATION:

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AU 2004262485	A1	20050217	AU 2004-262485	20040729
CA 2533976	A1	20050217	CA 2004-2533976	20040729
EP 1648445	A1	20060426	EP 2004-763610	20040729

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 IE, SI, LT, LV, FI, RO, CY, TR, BG, CZ, EE, HU, PL, SK
 CN 1832740 A 20060913 CN 2004-80022472 20040729
 BR 2004013110 A 20061003 BR 2004-13110 20040729
 JP 2007500165 T 20070111 JP 2006-521529 20040729
 MX 2006PA01159 A 20060424 MX 2006-PA1159 20060127
 NO 2006000865 A 20060222 NO 2006-865 20060222
 US 20070032520 A1 20070208 US 2006-566094 20061003
 PRIORITY APPLN. INFO.: ES 2003-1805 A 20030730
 OTHER SOURCE(S): CASREACT 142:240310; MARPAT 142:240310
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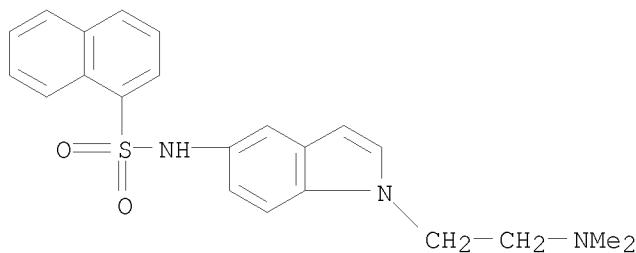


AB Title compds. I [R1 = NR8R9 radical or (un)saturated-(un)substituted cycloaliph. radical optionally containing at least one heteroatom; R2-4,6-7 independently = H, NO₂, alkoxy, CN, etc.; R5 = H or (un)saturated alkyl optionally at least monosubstituted; R8 or R9 independently = H or (un)saturated alkyl optionally at least monosubstituted with provisions; or R8 and R9 together with the bridging N atom form a (un)saturated-(un)substituted heterocyclic ring; A = (un)substituted mono or polycyclic aromatic ring; n = 0-4] and their pharmaceutically acceptable salts are prepared and disclosed as 5-HT₆ modulators. Thus, e.g., II, was prepared via reaction of naphthalene-2-sulfonyl chloride with 5-amino-1-(2-dimethylaminoethyl)-1H-indole. Selected data from 5-HT₆ receptor binding studies revealed Ki values (nM) ranging from 1.89-112.4.

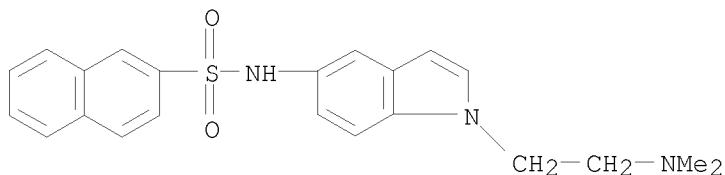
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 844832-10-6P

RL: PAC (Pharmacological activity); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)
 (drug candidate; preparation of indol-5-ylsulfonamide derivs. as 5-HT₆ receptor modulators)

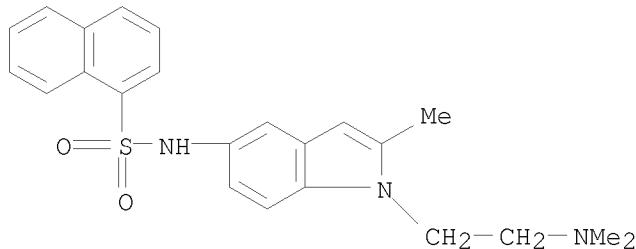
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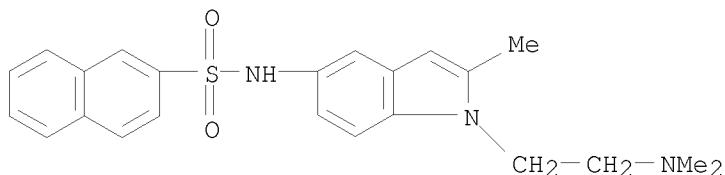
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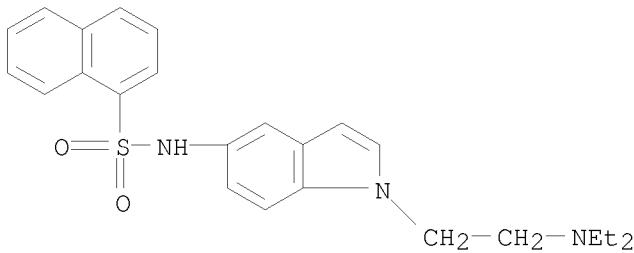
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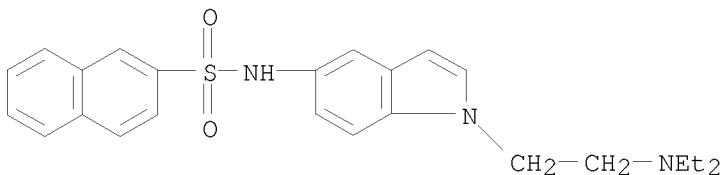


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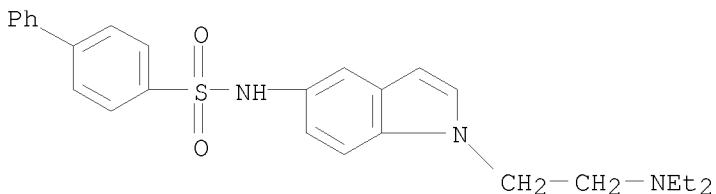
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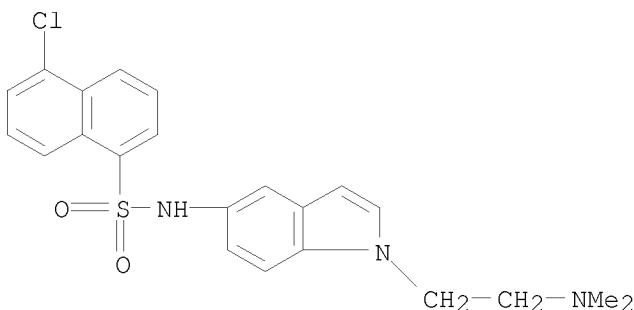
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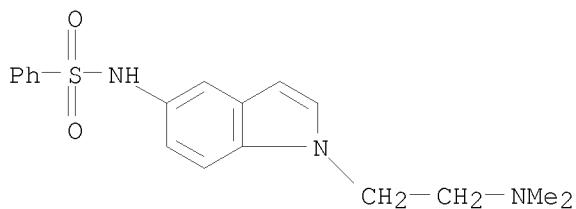
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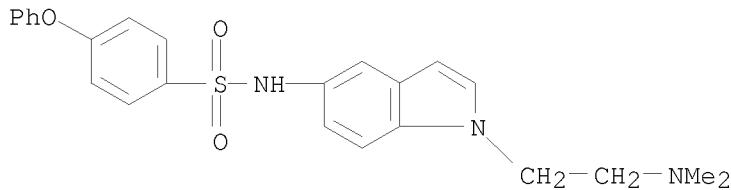


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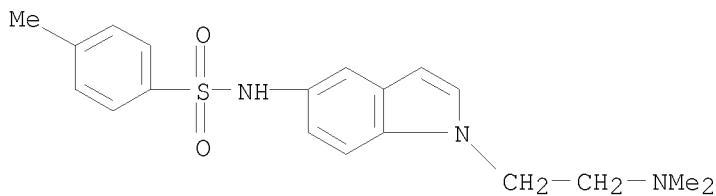
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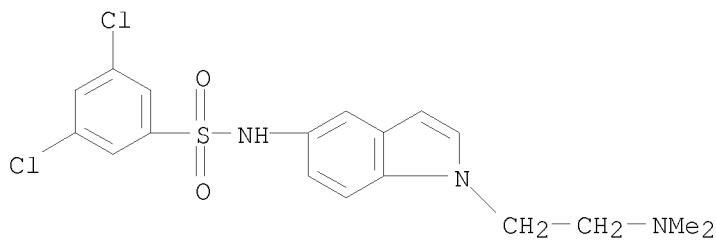
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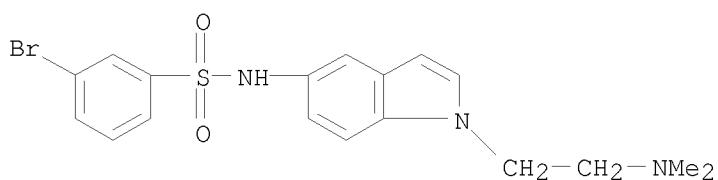
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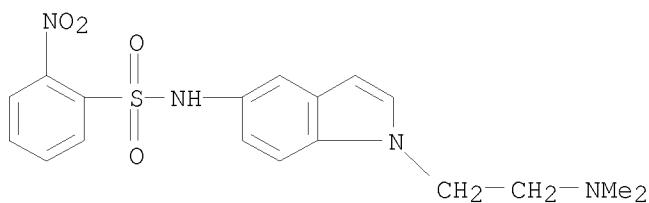
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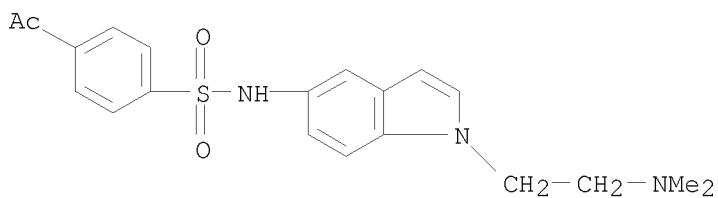
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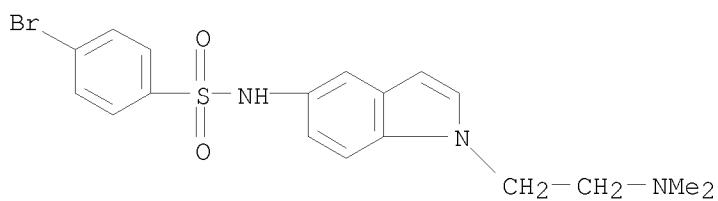
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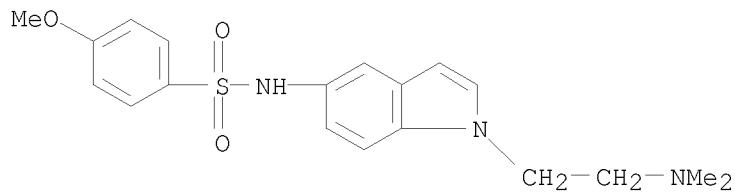
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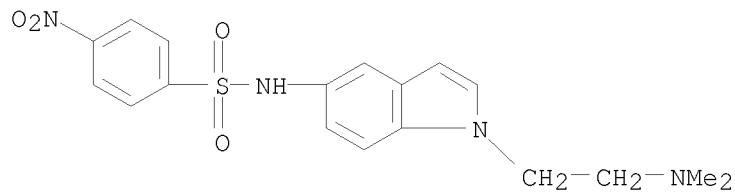


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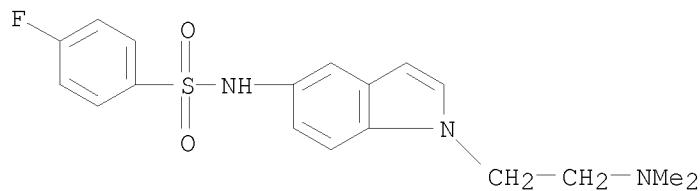
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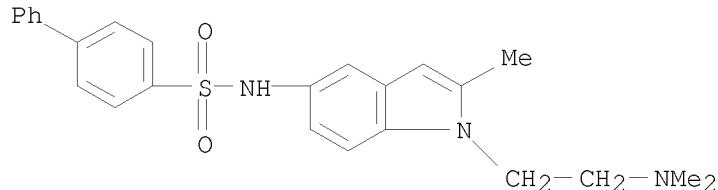
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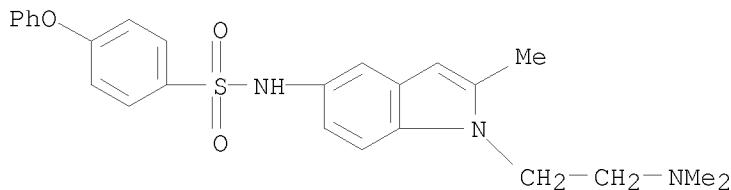
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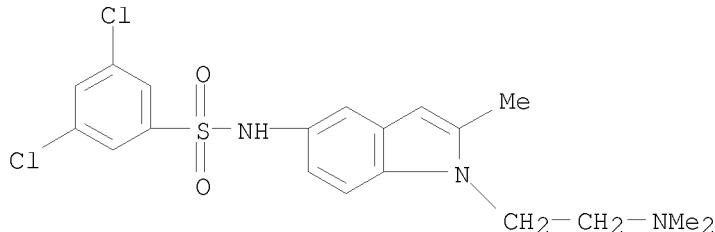
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RN 844832-10-6 CAPLUS

CN Benzenesulfonamide, 3,5-dichloro-N-[1-[2-(dimethylamino)ethyl]-2-methyl-1H-indol-5-yl]- (CA INDEX NAME)



REFERENCE COUNT: 5 THERE ARE 5 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

=> FILE REGISTRY

COST IN U.S. DOLLARS	SINCE FILE ENTRY	TOTAL SESSION
FULL ESTIMATED COST	37.50	216.99
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 DICTIONARY FILE UPDATES: 3 JUN 2008 HIGHEST RN 1025148-33-7

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<http://www.cas.org/support/stngen/stndoc/properties.html>

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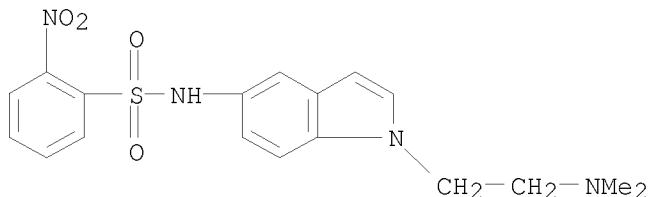
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L5 ANSWER 1 OF 1 REGISTRY COPYRIGHT 2008 ACS on STN
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(CA INDEX NAME)
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DT.CA CAplus document type: Patent
RL.P Roles from patents: BIOL (Biological study); PREP (Preparation); USES
(Uses)



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

3 REFERENCES IN FILE CA (1907 TO DATE)
3 REFERENCES IN FILE CAPLUS (1907 TO DATE)

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STRUCTURE FILE UPDATES: 3 JUN 2008 HIGHEST RN 1025148-33-7
DICTIONARY FILE UPDATES: 3 JUN 2008 HIGHEST RN 1025148-33-7

New CAS Information Use Policies, enter HELP USAGETERMS for details.

TSCA INFORMATION NOW CURRENT THROUGH January 9, 2008.

Please note that search-term pricing does apply when conducting SmartSELECT searches.

REGISTRY includes numerically searchable data for experimental and predicted properties as well as tags indicating availability of experimental property data in the original document. For information on property searching in REGISTRY, refer to:

<http://www.cas.org/support/stngen/stndoc/properties.html>

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COST IN U.S. DOLLARS	SINCE FILE ENTRY	TOTAL SESSION	
FULL ESTIMATED COST	0.46	219.91	
DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)	SINCE FILE ENTRY	TOTAL SESSION	
CA SUBSCRIBER PRICE	0.00	-4.80	

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FILE LAST UPDATED: 3 Jun 2008 (20080603/ED)

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=> s 15
L7 3 L5

=> d ibib abs 17 1-3

L7 ANSWER 1 OF 3 CAPLUS COPYRIGHT 2008 ACS on STN
ACCESSION NUMBER: 2005:136598 CAPLUS
DOCUMENT NUMBER: 142:240323
TITLE: Active substance combination comprising a compound with NPY receptor affinity and a compound with 5-HT6 receptor affinity
INVENTOR(S): Torrens Jover, Antoni; Mas Prio, Josep; Dordal Zueras, Alberto; Codony Soler, Xavier; Merce Vidal, Ramon; Aurelio Castrillo Perez, Jose; Frigola Constansa, Jordi; Buschmann, Helmut-Heinrich
PATENT ASSIGNEE(S): Laboratorios del Esteve S. A., Spain
SOURCE: PCT Int. Appl., 427 pp.
CODEN: PIXXD2
DOCUMENT TYPE: Patent
LANGUAGE: English
FAMILY ACC. NUM. COUNT: 1
PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
WO 2005014045	A1	20050217	WO 2004-EP8514	20040729
W: AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BW, BY, BZ, CA, CH, CN, CO, CR, CU, CZ, DE, DK, DM, DZ, EC, EE, EG, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NA, NI, NO, NZ, OM, PG, PH, PL, PT, RO, RU, SC, SD, SE, SG, SK, SL, SY, TJ, TM, TN, TR, TT, TZ, UA, UG, US, UZ, VC, VN, YU, ZA, ZM, ZW				
RW: BW, GH, GM, KE, LS, MW, MZ, NA, SD, SL, SZ, TZ, UG, ZM, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM, AT, BE, BG, CH, CY, CZ, DE, DK, EE, ES, FI, FR, GB, GR, HU, IE, IT, LU, MC, NL, PL, PT, RO, SE, SI, SK, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE, SN, TD, TG				
ES 2228268	A1	20050401	ES 2003-1815	20030730
ES 2228268	B1	20060701		
AU 2004262488	A1	20050217	AU 2004-262488	20040729
CA 2534099	A1	20050217	CA 2004-2534099	20040729
EP 1660131	A1	20060531	EP 2004-741321	20040729
R: AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, MC, PT, IE, SI, LT, LV, FI, RO, CY, TR, BG, CZ, EE, HU, PL, SK				
MX 2006PA01230	A	20060515	MX 2006-PA1230	20060130
US 20070009597	A1	20070111	US 2006-566402	20060705
PRIORITY APPLN. INFO.:			ES 2003-1815	A 20030730
			WO 2004-EP8514	W 20040729
OTHER SOURCE(S):	CASREACT 142:240323; MARPAT 142:240323			
GI				

* STRUCTURE DIAGRAM TOO LARGE FOR DISPLAY - AVAILABLE VIA OFFLINE PRINT *

AB The present invention relates to an active substance combination comprising at least one compound I [R1-R4 = H, halo, alkyl, etc.; R5 = H, alkyl, (un)saturated cycloalkyl; R6-R9 = H, alkyl, (un)saturated cycloalkyl, etc.]; A = CHR18, CHR18CH2; B = alkyl, (un)saturated cycloalkyl, etc.; R10 = H, alkyl, (un)saturated cycloalkyl, etc.; R11 = alkyl, (un)saturated cycloalkyl, etc.; NR10R11 = (un)saturated heterocyclyl; R18 = H, alkyl, (un)saturated cycloalkyl, etc.] with neuropeptide Y-receptor affinity, preferably

neuropeptide Y5-receptor affinity, and at least one compound with 5-HT6 receptor affinity (such as II [R1 = H, alkyl, Ph, CH2PH; R2 = NR4R5, (un)saturated (hetero)cycloalkyl, etc.; R3 = H, alkyl; R4, R5 = H, alkyl; or NR4R5 = (un)saturated heterocyclyl; A = (un)substituted (hetero)aryl; n = 0-4]), a medicament comprising said active substance combination, and the use of said active substance combination for the manufacture of a medicament. Synthesis of amides I and sulfonamides such as II is described in examples. E.g., a multi-step synthesis of III.HCl, starting from 1-(tert-butoxycarbonyl)-4-piperidinone and Me anthranilate, was given. The amides I and sulfonamides such as II were tested against neuropeptide Y5 and 5-HT6 binding (data given for representative compds.).

REFERENCE COUNT: 4 THERE ARE 4 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L7 ANSWER 2 OF 3 CAPLUS COPYRIGHT 2008 ACS on STN
 ACCESSION NUMBER: 2005:136568 CAPLUS
 DOCUMENT NUMBER: 142:240322
 TITLE: Active substance combination comprising a compound with NPY receptor affinity and a compound with 5-HT6 receptor affinity
 INVENTOR(S): Torrens Jover, Antoni; Mas Prio, Josep; Dordal Zueras, Alberto; Codony Soler, Xavier; Merce Vidal, Ramon; Aurelio Castrillo Perez, Jose; Frigola Constansa, Jordi; Buschmann, Helmut-Heinrich
 PATENT ASSIGNEE(S): Laboratorios del Esteve S. A., Spain
 SOURCE: PCT Int. Appl., 451 pp.
 CODEN: PIXXD2
 DOCUMENT TYPE: Patent
 LANGUAGE: English
 FAMILY ACC. NUM. COUNT: 1
 PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
WO 2005014000	A1	20050217	WO 2004-EP8515	20040729
W: AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BW, BY, BZ, CA, CH, CN, CO, CR, CU, CZ, DE, DK, DM, DZ, EC, EE, EG, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NA, NI, NO, NZ, OM, PG, PH, PL, PT, RO, RU, SC, SD, SE, SG, SK, SL, SY, TJ, TM, TN, TR, TT, TZ, UA, UG, US, UZ, VC, VN, YU, ZA, ZM, ZW				
RW: BW, GH, GM, KE, LS, MW, MZ, NA, SD, SL, SZ, TZ, UG, ZM, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM, AT, BE, BG, CH, CY, CZ, DE, DK, EE, ES, FI, FR, GB, GR, HU, IE, IT, LU, MC, NL, PL, PT, RO, SE, SI, SK, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE, SN, TD, TG				
ES 2228267	A1	20050401	ES 2003-1814	20030730
ES 2228267	B1	20060701		
AU 2004262489	A1	20050217	AU 2004-262489	20040729
CA 2534100	A1	20050217	CA 2004-2534100	20040729
EP 1648468	A1	20060426	EP 2004-763612	20040729
R: AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, MC, PT, IE, SI, LT, LV, FI, RO, CY, TR, BG, CZ, EE, HU, PL, SK				
MX 2006PA01232	A	20060515	MX 2006-PA1232	20060130
US 20070059364	A1	20070315	US 2006-566100	20061026
PRIORITY APPLN. INFO.:			ES 2003-1814	A 20030730
			WO 2004-EP8515	W 20040729

OTHER SOURCE(S): MARPAT 142:240322
 GI

* STRUCTURE DIAGRAM TOO LARGE FOR DISPLAY - AVAILABLE VIA OFFLINE PRINT *

AB The present invention relates to an active substance combination comprising at least one compound I [R1-R4 = H, halo, alkyl, etc.; R5 = H, alkyl, (un)saturated (hetero)cycloalkyl; R6-R9 = H, alkyl, (un)saturated (hetero)cycloalkyl, etc.; A = CHR18, CHR18CH2; R10 = H, alkyl, (un)saturated cycloalkyl, etc.; R11 = alkyl, (un)saturated cycloalkyl, etc.; NR10R11 = (un)saturated heterocycl; R18 = H, alkyl, (un)saturated cycloalkyl, etc.]

with

neuropeptide Y-receptor affinity, preferably neuropeptide Y5-receptor affinity, and at least one compound with 5-HT6 receptor affinity (such as II [R1 = H, alkyl, Ph, CH2PH; R2 = NR4R5, (un)saturated (hetero)cycloalkyl, etc.; R3 = H, alkyl; R4, R5 = H, alkyl; or NR4R5 = (un)saturated heterocycl; A = (un)substituted (hetero)aryl; n = 0-4]), a medicament comprising said active substance combination, and the use of said active substance combination for the manufacture of a medicament. Synthesis of amides I and sulfonamides such as II is described in examples. Thus, reacting 6-chloro-1-(4-piperidinyl)-1,4-dihydro-2H-3,1-benzoxazinone hydrochloride with 2-(2-chloroacetamide)-2',5-dichlorobenzophenone in the presence of K2CO3 in DMF followed by treating of the free base with HCl/EtOH afforded 61% III.HCl. The amides I and sulfonamides such as II were tested against neuropeptide Y5 and 5-HT6 binding (data given for representative compds.).

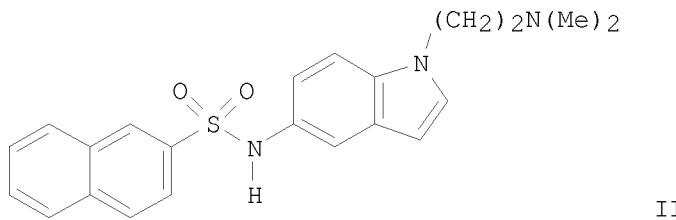
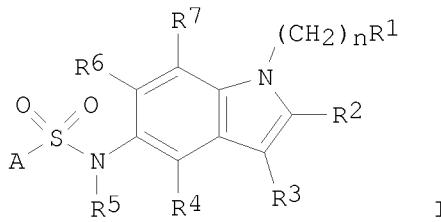
REFERENCE COUNT: 5 THERE ARE 5 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L7 ANSWER 3 OF 3 CAPLUS COPYRIGHT 2008 ACS on STN
ACCESSION NUMBER: 2005:136549 CAPLUS
DOCUMENT NUMBER: 142:240310
TITLE: Preparation of indol-5-yl sulfonamide derivatives and their use as 5-HT6 modulators
INVENTOR(S): Merce Vidal, Ramon; Codony Soler, Xavier; Dordal Zueras, Alberto
PATENT ASSIGNEE(S): Laboratorios del Esteve S. A., Spain
SOURCE: PCT Int. Appl., 123 pp.
CODEN: PIXXD2
DOCUMENT TYPE: Patent
LANGUAGE: English
FAMILY ACC. NUM. COUNT: 1
PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
WO 2005013977	A1	20050217	WO 2004-EP8511	20040729
W: AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BW, BY, BZ, CA, CH, CN, CO, CR, CU, CZ, DE, DK, DM, DZ, EC, EE, EG, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NA, NI, NO, NZ, OM, PG, PH, PL, PT, RO, RU, SC, SD, SE, SG, SK, SL, SY, TJ, TM, TN, TR, TT, TZ, UA, UG, US, UZ, VC, VN, YU, ZA, ZM, ZW				
RW: BW, GH, GM, KE, LS, MW, MZ, NA, SD, SL, SZ, TZ, UG, ZM, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM, AT, BE, BG, CH, CY, CZ, DE, DK, EE, ES, FI, FR, GB, GR, HU, IE, IT, LU, MC, NL, PL, PT, RO, SE, SI, SK, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE, SN, TD, TG				
ES 2222827	A1	20050201	ES 2003-1805	20030730
ES 2222827	B1	20060301		
AU 2004262485	A1	20050217	AU 2004-262485	20040729
CA 2533976	A1	20050217	CA 2004-2533976	20040729
EP 1648445	A1	20060426	EP 2004-763610	20040729
R: AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, MC, PT, IE, SI, LT, LV, FI, RO, CY, TR, BG, CZ, EE, HU, PL, SK				

CN 1832740	A	20060913	CN 2004-80022472	20040729
BR 2004013110	A	20061003	BR 2004-13110	20040729
JP 2007500165	T	20070111	JP 2006-521529	20040729
MX 2006PA01159	A	20060424	MX 2006-PA1159	20060127
NO 2006000865	A	20060222	NO 2006-865	20060222
US 20070032520	A1	20070208	US 2006-566094	20061003
PRIORITY APPLN. INFO.:			ES 2003-1805	A 20030730
			WO 2004-EP8511	W 20040729

OTHER SOURCE(S): CASREACT 142:240310; MARPAT 142:240310
GI



AB Title compds. I [R1 = NR8R9 radical or (un)saturated-(un)substituted cycloaliph. radical optionally containing at least one heteroatom; R2-4,6-7 independently = H, NO₂, alkoxy, CN, etc.; R5 = H or (un)saturated alkyl optionally at least monosubstituted; R8 or R9 independently = H or (un)saturated alkyl optionally at least monosubstituted with provisions; or R8 and R9 together with the bridging N atom form a (un)saturated-(un)substituted heterocyclic ring; A = (un)substituted mono or polycyclic aromatic ring; n = 0-4] and their pharmaceutically acceptable salts are prepared and disclosed as 5-HT₆ modulators. Thus, e.g., II, was prepared via reaction of naphthalene-2-sulfonyl chloride with 5-amino-1-(2-dimethylaminoethyl)-1H-indole. Selected data from 5-HT₆ receptor binding studies revealed Ki values (nM) ranging from 1.89-112.4.

REFERENCE COUNT: 5 THERE ARE 5 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

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ALL L# QUERIES AND ANSWER SETS ARE DELETED AT LOGOFF

LOGOFF? (Y)/N/HOLD:y

COST IN U.S. DOLLARS	SINCE FILE ENTRY	TOTAL SESSION
FULL ESTIMATED COST	9.21	229.12
DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)	SINCE FILE ENTRY	TOTAL SESSION

CA SUBSCRIBER PRICE

-2.40

-7.20

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